

Simulation of crashworthiness problems with improved implicit time integration methods for non-linear dynamics.

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- Industrial context:
 - Structures must be able to resist to crash situations
 - Numerical simulations is a key to design structures
 - Efficient time integration in the non-linear range is needed
- Goal:
 - Numerical simulation of
 blade off and wind-milling in
 a turboengine
 - Example from SNECMA





- 1. Scientific motivations
- 2. Consistent scheme in the non-linear range
- 3. Combined implicit/explicit algorithm
- 4. Complex numerical examples
- 5. Conclusions & perspectives



- Scientific context:
 - Solids mechanics
 - Large displacements
 - Large deformations
 - Non-linear mechanics
- Spatial discretization into finite elements:
 - Balance equation $M \ddot{\vec{x}} + \vec{F}^{\text{int}} = \vec{F}^{\text{ext}}$
 - Internal forces formulation

 $\vec{F}^{\text{int}} = \int_{V_0} \Sigma \mathbf{f}^T \vec{D} J dV_0$

Σ: Cauchy stress; **F**: deformation gradient; $\mathbf{f} = \mathbf{F}^{-1}$; $\vec{D} = \partial \varphi / \partial \vec{x}_0$: derivative of the shape function; *J*: Jacobian



1. Scientific motivations Dynamics simulations

- Temporal integration of the balance equation
- 2 methods: - Explicit method $\begin{vmatrix} \vec{x}_n \\ \vec{x}_n \\ \vec{x}_n \end{vmatrix}$ approximation $\Rightarrow \vec{x}_{n+1} = \mathbf{M}^{-1} \left(\vec{F}_n^{\text{ext}} - \vec{F}_n^{\text{int}} \right) \xrightarrow{\text{deduction}} \vec{x}_{n+1}, \dot{\vec{x}}_{n+1}$ Very fast dynamics • Non iterative • Limited needs in memory • Conditionally stable (small time step) $- \text{ Implicit method } \begin{vmatrix} \vec{x}_n \\ \dot{\vec{x}}_n \\ \vdots \\ \vec{x}_n \end{vmatrix} \xrightarrow{\text{extrapolation}} \begin{cases} \vec{x}_{n+1} \\ \dot{\vec{x}}_{n+1} \\ \vdots \\ \vec{x}_{n+1} \end{cases} \xleftarrow{\text{iterations}} \begin{cases} \mathbf{M} \ddot{\vec{x}} + \vec{F}^{\text{int}} = \vec{F}^{\text{ext}} \\ \vec{x}_{n+1} = f(\vec{x}_n, \dot{\vec{x}}_n, \dot{\vec{x}}_{n+1}, \ddot{\vec{x}}_n, \ddot{\vec{x}}_{n+1}) \\ \dot{\vec{x}}_{n+1} = f(\vec{x}_n, \vec{x}_n, \vec{x}_{n+1}, \ddot{\vec{x}}_n, \ddot{\vec{x}}_{n+1}) \end{cases}$ Slower dynamics • Iterative • More needs in memory • Unconditionally stable (large time step)



- If wave propagation effects are negligible
 - ---- Implicit schemes are more suitable
 - Sheet metal forming (springback, superplastic forming, ...)
 - Crashworthiness simulations (car crash, blade loss, shock absorber, ...)
- Nowadays, people choose explicit scheme mainly because of difficulties linked to implicit scheme:
 - Lack of smoothness (contact, elasto-plasticity, ...)
 - ---- convergence can be difficult
 - Lack of available methods (commercial codes)
- Little room for improvement in explicit methods
- Complex problems can take advantage of combining explicit and implicit algorithms



1. Scientific motivations Conservation laws

- Conservation of linear momentum (Newton's law)
 - Continuous dynamics
 - Time discretization

- Conservation of angular momentum
 - Continuous dynamics
 - Time discretization

$$\frac{\partial \vec{x} \wedge \mathbf{M} \vec{x}}{\partial t} = \vec{x} \wedge \vec{F}^{\text{ext}}$$

$$\sum_{nodes} \sum_{nodes} \vec{x}_{n+1} \wedge \mathbf{M} \dot{\vec{x}}_{n+1} - \vec{x}_n \wedge \mathbf{M} \dot{\vec{x}}_n = \Delta t \sum_{nodes} \vec{x}_{n+1/2} \wedge \vec{F}_{n+1/2}^{\text{ext}}$$

$$\bigotimes_{nodes} \sum_{nodes} \vec{x}_{n+1/2} \wedge \vec{F}_{n+1/2}^{\text{int}} = 0$$
Wint is to all

Conservation of energy

W^{int}: internal energy; W^{ext}: external energy;

D^{int}: dissipation (plasticity ...)

0

- Continuous dynamics $\frac{\partial}{\partial t}K + \frac{\partial}{\partial t}W^{\text{int}} = \frac{\partial}{\partial t}W^{\text{ext}} \dot{D}^{\text{int}}$
- Time discretization $W_{n+1}^{\text{int}} W_n^{\text{int}} + \Delta D^{\text{int}} = \sum_{nodes} \vec{F}_{n+1/2}^{\text{int}} \cdot [\vec{x}_{n+1} \vec{x}_n] \&$

$$\sum_{nodes} \frac{1}{2} \mathbf{M} \dot{\vec{x}}_{n+1} \bullet \dot{\vec{x}}_{n+1} - \frac{1}{2} \mathbf{M} \dot{\vec{x}}_n \bullet \dot{\vec{x}}_n + W_{n+1}^{\text{int}} - W_n^{\text{int}} + \Delta D^{\text{int}} = \sum_{nodes} \vec{F}_{n+1/2}^{\text{ext}} \bullet [\vec{x}_{n+1} - \vec{x}_n]$$



1. Scientific motivations Explicit algorithms

Central difference (no numerical dissipation)

 $\dot{\vec{x}}_{n+1/2} = \dot{\vec{x}}_{n-1/2} + \Delta t \quad \ddot{\vec{x}}_n$ $\vec{x}_{n+1} = \vec{x}_n + \Delta t \quad \dot{\vec{x}}_{n+1/2}$ $\ddot{\vec{x}}_{n+1} = M^{-1} \Big[\vec{F}_{n+1}^{\text{ext}} - \vec{F}_{n+1}^{\text{int}} \Big]$

- Hulbert & Chung (numerical dissipation) [CMAME, 1996] $\begin{bmatrix} 1-\alpha_{M} \end{bmatrix} \ddot{\vec{x}}_{n+1} = M^{-1} \begin{bmatrix} \vec{F}_{n}^{\text{ext}} - \vec{F}_{n}^{\text{int}} \end{bmatrix} - \alpha_{M} \ddot{\vec{x}}_{n}$ $\vec{x}_{n+1} = \vec{x}_{n} + \Delta t \, \dot{\vec{x}}_{n} + \Delta t^{2} \begin{bmatrix} \frac{1}{2} - \beta \end{bmatrix} \ddot{\vec{x}}_{n} + \Delta t^{2} \beta \, \ddot{\vec{x}}_{n+1}$ $\dot{\vec{x}}_{n+1} = \dot{\vec{x}}_{n} + \Delta t [1-\gamma] \ddot{\vec{x}}_{n} + \Delta t \gamma \, \ddot{\vec{x}}_{n+1}$
- Small time steps conservation conditions are approximated
- Numerical oscillations may cause spurious plasticity



1. Scientific motivations Implicit algorithms

- α-generalized family (Chung & Hulbert [JAM, 1993])
 - Newmark relations: $\begin{cases} \ddot{x}_{n+1} = \frac{1}{\beta \Delta t^2} \left[\vec{x}_{n+1} \vec{x}_n \Delta t \, \dot{\vec{x}}_n \left[\frac{1}{2} \beta \right] \Delta t^2 \, \ddot{\vec{x}}_n \right] \\ \dot{\vec{x}}_{n+1} = \frac{\gamma}{|\vec{x}_{n+1} \vec{x}_n|} \left[\vec{x}_{n+1} \vec{x}_n + \left[\frac{\beta}{|\vec{x}_n|} \frac{1}{2} \right] \Delta t^2 \, \ddot{\vec{x}}_n \right] \end{cases}$

$$\frac{1-\alpha_M}{1-\alpha_F}M\ddot{\vec{x}}_{n+1} + \frac{\alpha_M}{1-\alpha_F}M\ddot{\vec{x}}_n + \left[\vec{F}_{n+1}^{\text{int}} - \vec{F}_{n+1}^{\text{ext}}\right] + \frac{\alpha_F}{1-\alpha_F}\left[\vec{F}_n^{\text{int}} - \vec{F}_n^{\text{ext}}\right] = 0$$

 $- \alpha_{\rm M} = 0$ and $\alpha_{\rm F} = 0$ (no numerical dissipation)

- Linear range: consistency (i.e. physical results) demonstrated
- Non-linear range with small time steps: consistency verified
- Non-linear range with large time steps: total energy conserved but without consistency (e.g. plastic dissipation greater than the total energy, work of the normal contact forces > 0, ...)
- $\alpha_M \neq 0$ and/or $\alpha_F \neq 0$ (numerical dissipation)
 - Numerical dissipation is proved to be positive only in the linear range

1. Scientific motivations Numerical example: mass/spring-system

- Example: Mass/spring system (2D) with an initial velocity perpendicular to the spring (Armero & Romero [CMAME, 1999])
 - Newmark implicit scheme (no numerical damping)





Explicit method: ∆tcrit ~ 0.72s; 1 revolution ~ 4s

Chung-Hulbert implicit scheme (numerical damping)



1. Scientific motivations Numerical example: mass/spring-system

- Same mass/spring system with a central difference scheme
 - $\Delta t_{crit} \sim 0.72$ s computed from the maximal pulsation of the system (2 degrees of freedom)



2. Consistent scheme in the non linear range Principle

- Consistent implicit algorithms in the non-linear range:
 - The Energy Momentum Conserving Algorithm or EMCA (Simo et al. [ZAMP 92], Gonzalez & Simo [CMAME 96]):
 - Conservation of the linear momentum
 - Conservation of the angular momentum
 - Conservation of the energy (no numerical dissipation)
 - The Energy Dissipative Momentum Conserving algorithm or EDMC (Armero & Romero [CMAME, 2001]):
 - Conservation of the linear momentum
 - Conservation of the angular momentum
 - Numerical dissipation of the energy is proved to be positive

2. Consistent scheme in the non linear range Principle

- Based on the mid-point scheme (Simo et al. [ZAMP, 1992])
 - Relations displacements
 /velocities/accelerations
 - Balance equation

$$\begin{cases} \frac{\ddot{x}_{n+1} + \ddot{x}_n}{2} = \frac{\dot{x}_{n+1} - \dot{x}_n}{\Delta t} \\ \frac{\dot{x}_{n+1} + \dot{x}_n}{2} \left(+ \dot{x}_{n+1}^{\text{diss}} \right) = \frac{\vec{x}_{n+1} - \vec{x}_n}{\Delta t} \\ \frac{\ddot{x}_{n+1} + \ddot{x}_n}{2} = \vec{F}_{n+1/2}^{\text{ext}} - \vec{F}_{n+1/2}^{\text{int}} \left(- \vec{F}_{n+1/2}^{\text{diss}} \right) \end{cases}$$

– EMCA:

- With $\vec{F}_{n+1/2}^{\text{int}} \neq \int \Sigma^{n+1/2} \mathbf{f}_0^{n+1/2^T} \vec{D} J_0^{n+1/2} dV_0$ and $\vec{F}_{n+1/2}^{\text{ext}}$ designed to verify conserving equations
- No dissipation forces and no dissipation velocities EDMC:
 - Same internal and external forces as in the EMCA
 - With $\vec{F}_{n+1/2}^{\text{diss}}$ and $\vec{x}_{n+1}^{\text{diss}}$ designed to achieve positive numerical dissipation without spectral bifurcation

2. Consistent scheme in the non linear range Dissipation property

- Comparison of the spectral radius
 - Integration of a linear oscillator: \vec{x}_n

$$S_{n+1} = \exp\left(\frac{\ln \rho^2}{\Delta t}t_{n+1}\right)\cos\left(\omega t_{n+1}\right)$$

 ρ : spectral radius; ω : pulsation

Low numerical dissipation

High numerical dissipation



2. Consistent scheme in the non linear range The mass/spring system

- Forces of the spring for any potential V
 - <u>Without</u> numerical dissipation \vec{F}_{n}^{\dagger} (EMCA) (Gonzalez & Simo [CMAME, 1996])

$$\lim_{n+1/2} = \frac{V(l_{n+1}) - V(l_n)}{l_{n+1}^2 - l_n^2} \left[\vec{x}_{n+1} + \vec{x}_n \right]$$



- The consistency of the EMCA solution does not depend on Δt
- The Newmark solution does-not conserve the angular momentum

2. Consistent scheme in the non linear range The mass/spring system

 $\dot{\vec{x}}_{n+1}^{\text{diss}} = \chi \frac{\|\vec{x}_{n+1}\| - \|\vec{x}_n\|}{\|\dot{\vec{x}}_{n+1}\| + \|\dot{\vec{x}}_n\|} \frac{[\vec{x}_{n+1} + \vec{x}_n]}{2}$

– <u>With</u> numerical dissipation (EDMC 1st order) with dissipation parameter $0 < \chi < 1$ (Armero&Romero [CMAME, 2001]), here $\chi = 0.111$



- Only EDMC solution preserves the driving motion:
 - The length tends towards the equilibrium length
 - Conservation of the angular momentum is achieved

EXAMPLE 2. Consistent scheme in the non linear range Formulations in the literature: hyperelasticity

- Hyperelastic material (stress derived from a potential *V*):
 - Saint Venant-Kirchhoff hyperelastic model (Simo et al. [ZAMP, 1992])
 - General formulation for hyperelasticity (Gonzalez [CMAME, 2000]):

$$\vec{F}_{n+1/2}^{\text{int}} = \int_{V_0} \frac{\mathbf{F}_0^n + \mathbf{F}_0^{n+1}}{2} \begin{bmatrix} \frac{\partial V}{\partial \mathbf{GL}} \left(\frac{\mathbf{GL}_0^n + \mathbf{GL}_0^{n+1}}{2} \right) + \tilde{f(V^{n+1}, V^n)} \end{bmatrix} \vec{D}dV_0$$

F: deformation gradient

GL: Green-Lagrange strain

V: potential

$$\vec{D} = \partial \varphi / \partial \vec{x}_0$$

 φ : shape functions

- Classical formulation: $\vec{F}^{\text{int}} = \int_{V_0} \mathbf{F} \frac{\partial V}{\partial \mathbf{GL}} \vec{D} dV_0$
- Hyperelasticity with elasto-plastic behavior: energy dissipation of the algorithm corresponds to the internal dissipation of the material (Meng & Laursen [CMAME, 2001])

2. Consistent scheme in the non linear range Formulations in the literature: contact

Description of the contact interaction:



- Computation of the classical contact force:
 - Penalty method $\vec{F}^{\text{cont}} = -k_N g \vec{n}$ k_N : penalty
 - Augmented Lagrangian method $\vec{F}^{\text{cont}} = \Lambda^{(k)} \vec{n} k_N g \vec{n}$

Λ: Lagrangian

- Lagrangian method $\vec{F}^{\text{cont}} = \Lambda \vec{n}$

EXAMPLE 2. Consistent scheme in the non linear range Formulations in the literature: contact

- Penalty contact formulation (normal force proportional to the penetration "gap") (Armero & Petöcz [CMAME, 1998-1999]):
 - Computation of a dynamic gap for slave node \vec{x} projected on master surface $\vec{y}(u)$

 $g_{n+1}^{d} = g_{n}^{d} + \vec{n}_{n+1/2} \bullet \left[\vec{x}_{n+1} - \vec{x}_{n} - \vec{y}_{n+1}(u_{n+1/2}) + \vec{y}_{n}(u_{n+1/2}) \right]$

Normal forces derived from a potential V

$$\vec{F}_{n+1/2}^{\text{cont}} = \frac{V(g_{n+1}^{d}) - V(g_{n}^{d})}{g_{n+1}^{d} - g_{n}^{d}} \vec{n}_{n+1/2}$$

- Augmented Lagrangian and Lagrangian consistent contact formulation (Chawla & Laursen [IJNME, 1997-1998]):
 - Computation of a gap rate

$$\Delta t \ g_{n+1/2}^{\mathbf{r}} = \vec{n}_{n+1/2} \bullet \left[\vec{x}_{n+1} - \vec{x}_n - \vec{y}_{n+1}(u_{n+1/2}) + \vec{y}_n(u_{n+1/2}) \right]$$

2. Consistent scheme in the non linear range Developments for a hypoelastic model

- The EMCA or EDMC for hypoelastic constitutive model:
 - Valid for hypoelastic formulation of (visco) plasticity
 - Energy dissipation from the internal forces corresponds to the plastic dissipation
- Hypoelastic model:
 - stress obtained incrementally from a hardening law
 - no possible definition of an internal potential!
 - Idea: the internal forces are established to be consistent on a loading/unloading cycle



strain

2. Consistent scheme in the non linear range Developments for a hypoelastic model

- Incremental strain tensor:
 E: natural strain tensor; F: deformation gradient
- Elastic incremental stress:
 Σ: Cauchy stress; *H*: Hooke stress-strain tensor
- Plastic stress corrections:

(radial return mapping: Wilkins [MCP, 1964], Maenchen & Sack [MCP, 1964], Ponthot [IJP, 2002])

s^c: plastic corrections; σ^{vm}: yield stress; ε^p: equivalent plastic strain

Final Cauchy stresses:

(final rotation scheme: Nagtegaal & Veldpaus [NAFP, 1984], Ponthot [IJP, 2002]) **R**: rotation tensor;

- Classical forces formulation:
 f = E⁻¹: D derivative of the shape function:
 - $\mathbf{f} = \mathbf{F}^{-1}$; *D* derivative of the shape function; *J* : Jacobian

$$\mathbf{E}_{n}^{n+1} = \frac{1}{2} \ln \left(\mathbf{F}_{n}^{n+1} \mathbf{F}_{n}^{n+1} \right)$$

$$\Delta \Sigma_n^{n+1} = H : \mathbf{E}_n^{n+1}$$

 $s_n^{c^{n+1}} = f(\sigma^{vm}, \varepsilon^p)$

$$\Sigma^{n+1} = \mathbf{R}_n^{n+1} \left[\Sigma^n + \Delta \Sigma_n^{n+1} - \mathbf{s}^c \right] \mathbf{R}_n^{n+1} \mathbf{T}$$

$$\vec{F}_{n+1}^{\text{int}} = \int_{V_0} \Sigma^{n+1} \mathbf{f}_0^{n+1^T} \vec{D} J dV_0$$

2. Consistent scheme in the non linear range Developments for a hypoelastic model

- EMCA (<u>without</u> numerical dissipation):
 - Balance equation $M \frac{\ddot{\vec{x}}_{n+1} + \ddot{\vec{x}}_n}{2} = \vec{F}_{n+1/2}^{\text{ext}} \vec{F}_{n+1/2}^{\text{int}}$
 - New internal forces formulation:

 $\vec{F}_{n+1/2}^{\text{int}} = \frac{1}{4} \int_{V_0} \left[\mathbf{I} + \mathbf{F}_n^{n+1} \right] \left[\Sigma^n + \mathbf{C}^* \right] \mathbf{f}_0^{n^T} \vec{D} J^n + \left[\mathbf{I} + \mathbf{f}_n^{n+1} \right] \left[\Sigma^{n+1} + \mathbf{C}^{**} \right] \mathbf{f}_0^{n+1^T} \vec{D} J^{n+1} dV_0$

F: deformation gradient; **f**: inverse of **F**; *D* derivative of the shape function; *J* : Jacobian = det **F**; Σ : Cauchy stress

 Correction terms C* and C**: (second order correction in the plastic strain increment)

$$\mathbf{C}^{*} = \frac{\Delta D^{\text{int}} / J_0^n - \mathbf{GL}_n^{n+1} p^l}{\mathbf{GL}_n^{n+1} : \mathbf{GL}_n^{n+1}} \mathbf{GL}_n^{n+1}$$
$$\mathbf{C}^{**} = \frac{\Delta D^{\text{int}} / J_0^{n+1} - \mathbf{A}_n^{n+1} p^l}{\mathbf{A}_n^{n+1} : \mathbf{A}_n^{n+1}} \mathbf{A}_n^{n+1}$$

 ΔD^{int} : internal dissipation due to the plasticity; **A**: Almansi incremental strain tensor ($\mathbf{A} = \mathbf{A}^{pl} + \mathbf{A}^{el}$); **GL**: Green-Lagrange incremental strain ($\mathbf{GL} = \mathbf{GL}^{pl} + \mathbf{GL}^{el}$)

- Verification of the conservation laws $\begin{cases} \sum_{nodes} \vec{F}_{n+1/2}^{\text{int}} = 0 & \& & \sum_{nodes} \vec{x}_{n+1/2} \land \vec{F}_{n+1/2}^{\text{int}} = 0 \\ \Delta D^{\text{int}} = \int_{cycle nodes} \sum_{rodes} \vec{F}_{n+1/2}^{\text{int}} \bullet [\vec{x}_{n+1} - \vec{x}_n] \end{cases}$

2. Consistent scheme in the non linear range Developments for a hypoelastic model

- EDMC (1st order accurate with numerical dissipation):
 - Balance equation $M \frac{\ddot{\vec{x}}_{n+1} + \ddot{\vec{x}}_n}{2} = \vec{F}_{n+1/2}^{\text{ext}} \vec{F}_{n+1/2}^{\text{int}} \vec{F}_{n+1/2}^{\text{diss}}$
 - New dissipation forces formulation:

$$\vec{F}_{n+1/2}^{\text{diss}} = \frac{1}{4} \int_{V_0} \left[\mathbf{I} + \mathbf{F}_n^{n+1} \right] \mathbf{D}^* \mathbf{f}_0^{n^T} \vec{D} + \left[\mathbf{I} + \mathbf{f}_n^{n+1} \right] \mathbf{D}^{**} \mathbf{f}_0^{n+1^T} \vec{D} dV_0$$

– Dissipating terms D* and D**:

$$\begin{cases} \mathbf{D}^{*} = \frac{\chi_{2}^{\prime} \mathbf{E}^{\mathbf{e}_{n}^{n+1}} : H : \mathbf{E}^{\mathbf{e}_{n}^{n+1}} J_{0}^{n}}{\mathbf{GL}_{n}^{n+1}} : \mathbf{GL}_{n}^{n+1} \\ \mathbf{D}^{**} = \frac{\chi_{2}^{\prime} / \mathbf{E}^{\mathbf{e}_{n}^{n+1}} : H : \mathbf{E}^{\mathbf{e}_{n}^{n+1}} J_{0}^{n+1}}{\mathbf{A}_{n}^{n+1}} : \mathbf{A}_{n}^{n+1} \end{cases}$$

Verification of the conservation laws

$$\begin{cases} \sum_{nodes} \vec{F}_{n+1/2}^{\text{diss}} = 0 \quad \& \quad \sum_{nodes} \vec{x}_{n+1/2} \wedge \vec{F}_{n+1/2}^{\text{diss}} = 0 \\ \Delta D^{\text{num}} = \sum_{nodes} \vec{F}_{n+1/2}^{\text{diss}} \bullet [\vec{x}_{n+1} - \vec{x}_n] + \mathbf{M} \dot{\vec{x}}_{n+1/2}^{\text{diss}} \bullet [\dot{\vec{x}}_{n+1} - \dot{\vec{x}}_n] \end{cases}$$

 ΔD^{num} : numerical dissipation

2. Consistent scheme in the non linear range Numerical example: Taylor bar

- Impact of a cylinder :
 - Hypoelastic model
 - Elasto-plastic hardening law
 - Simulation during 80 µs



2. Consistent scheme in the non linear range Numerical example: Taylor bar

Simulation without numerical dissipation: final results



2. Consistent scheme in the non linear range Numerical example: Taylor bar

Simulations with numerical dissipation: final results

- Constant spectral radius at infinity pulsation = 0.7



2. Consistent scheme in the non linear range Numerical example: impact of two 2D-cylinders

- Impact of 2 cylinders (Meng&Laursen) :
 - Left one has a initial velocity (initial kinetic energy 14J)
 - Elasto perfectly plastic hypoelastic material
 - Simulation during 4s



2. Consistent scheme in the non linear range Numerical example: impact of two 2D-cylinders

- Results comparison at the end of the simulation
 - Newmark($\Delta t=1.875$ ms)

EMCA (with cor., $\Delta t=1.875$ ms)



2. Consistent scheme in the non linear range Numerical example: impact of two 2D-cylinders

• Results evolution comparison $\Delta t=1.875$ ms





2. Consistent scheme in the non linear range Numerical example: impact of two 3D-cylinders

- Impact of 2 hollow 3D-cylinders:
 - Right one has a initial velocity $(\dot{\vec{x}}_{0X} = 10\dot{\vec{x}}_{0Y})$
 - Elasto-plastic
 hypoelastic material
 (aluminum)
 - Simulation during 5ms
 - Use of numerical dissipation
 - Frictional contact

y

Х



Numerical example: impact of two 3D-cylinders

• Results comparison with a reference (EMCA; $\Delta t=0.5\mu s$):



– At the end of the simulation:





0.0025

Time (s)

y

0.005

Х

Extension to the use of incremental potential

 Hyperelastic material with use of the variational formulation of visco-plastic updates [Ortiz & Stainier, CMAME 1999]

 $\mathbf{S}^{n+1} = 2 \frac{\partial \Delta D_{\text{eff}}}{\partial \mathbf{C}^{n+1}} \left(\mathbf{C}_0^{n+1}, \mathbf{C}_0^n \right)$

C: right Cauchy-Green strain

S: second Piola-Kirchhoff stress

 $\Delta D_{\rm eff}$: incremental potential

Use of a similar form than the Gonzalez formulation
 [CMAME 2000] for an elastic model

$$\vec{F}_{n+1/2}^{\text{int}} = \int_{V_0} \frac{\mathbf{F}_0^n + \mathbf{F}_0^{n+1}}{2} \left[2 \frac{\partial \Delta D_{\text{eff}}}{\partial \mathbf{C}} \left(\frac{\mathbf{C}_0^n + \mathbf{C}_0^{n+1}}{2}, \mathbf{C}_0^n \right) + f \left(\Delta D_{\text{eff}} \left(\mathbf{C}_0^{n+1}, \mathbf{C}_0^n \right), \frac{\partial \Delta D_{\text{eff}}}{\partial \mathbf{C}} \left(\frac{\mathbf{C}_0^n + \mathbf{C}_0^{n+1}}{2}, \mathbf{C}_0^n \right) \right) \right] \vec{D} dV_0$$

F: deformation gradient; φ : shape functions; $\vec{D} = \partial \varphi / \partial \vec{x}_0$

2. Consistent scheme in the non linear range Simulation of a tumbling beam

0.000

0.0325

- Tumbling beam:
 - Initial symmetrical loads (t < 10s)
 - Elasto perfectly
 plastic hyperelastic
 material
 - Simulation during 150s
 - Use of the variational formulation of elastoplastic updates
 - Conserving algorithm



0.0650

0.0975

0.130

2. Consistent scheme in the non linear range Simulation of a tumbling beam

Time evolution of the results:





3. Combined implicit/explicit algorithm Automatic shift

- Shift from an implicit algorithm to an explicit one:
 - Evaluation of the ratio r*
 - Explicit time step size depends on the mesh
 - Ωb : stability limit;
 - ω max: maximal eigen pulsation
 - Implicit time step size depends on the integration error (Géradin)
 *e*_{int}: integration error; *Tol*: user tolerance

$$r^* = \frac{CPU \quad 1 \quad \text{implicit} \quad \text{step}}{CPU \quad 1 \quad \text{explicit} \quad \text{step}}$$
$$\Delta t_{\text{expl}} = \frac{\Omega_b}{\omega_{\text{max}}}$$

$$\begin{cases} e_{\text{int}} = \Delta t^{3} \ddot{x} \approx \frac{\Delta t_{\text{impl}}^{2} \sum_{i=nodes} \Delta |\vec{x}|_{i}}{e_{\text{ref}}} \\ \frac{\Delta t_{\text{impl}}^{\text{new}}}{\Delta t_{\text{impl}}^{\text{old}}} = \left[\frac{2e_{\text{int}}}{Tol}\right]^{1/2.5} \end{cases}$$

$$\Delta t_{\rm impl} < \frac{r^* \Delta t_{\rm expl}}{\mu}$$

- Shift criterion
 - μ : user security



- Shift from an explicit algorithm to an implicit one:
 - Evaluation of the ratio r*
 - Explicit time step size depends on the mesh
 Ωb: stability limit;
 ω max: maximal eigen pulsation
 - Implicit time step size
 interpolated form a acceleration
 difference
 Tol: user tolerance
 - Shift criterion
 - μ : user security

$$r^* = \frac{CPU \quad 1 \quad \text{implicit} \quad \text{step}}{CPU \quad 1 \quad \text{explicit} \quad \text{step}}$$
$$\Delta t_{\text{expl}} = \frac{\Omega_b}{\omega_{\text{max}}}$$

$$\Delta t_{\text{impl}} = \left(\frac{Tol \ e_{\text{ref}} \sqrt{\Delta t_{\text{expl}}}}{2\sum_{i=nodes} \Delta \left|\vec{x}\right|_{i}}\right)^{2/5}$$

$$\Delta t_{\rm impl} > \mu r^* \Delta t_{\rm expl}$$



Stabilization of the explicit solution:



- Dissipation of the numerical modes: spectral radius at bifurcation equal to zero.
- Consistent balance of the r* last explicit steps:

$$M \frac{\ddot{\vec{x}}_{n+r^*} + \ddot{\vec{x}}_n}{2} = \vec{F}_{n+r^*/2} - \vec{F}_{n+r^*/2} \left(-\vec{F}_{n+r^*/2} \right)$$

. 3. Combined implicit/explicit algorithm Numerical example: blade casing interaction

- Blade/casing interaction :
 - Rotation velocity3333rpm
 - Rotation center is moved during the first half revolution
 - EDMC-1 algorithm
 - Four revolutions simulation



3. Combined implicit/explicit algorithm Numerical example: blade casing interaction

• Final results comparison:



4. Complex numerical examples Blade off simulation

Numerical simulation of a blade loss in an aero engine





4. Complex numerical examples Blade off simulation

- Blade off :
 - Rotation velocity5000rpm
 - EDMC algorithm
 - 29000 dof's
 - One revolution simulation
 - 9000 time steps
 - 50000 iterations (only
 9000 with stiffness
 matrix updating)





4. Complex numerical examples Blade off simulation

Final results comparison:





• CPU time comparison before and after code optimization: Before optimization







4. Complex numerical examples Dynamic buckling of square aluminum tubes

- Absorption of 600J with different impact velocities :
 - EDMC algorithm
 - 16000 dof's / 2640 elements
 - Initial asymmetry
 - Comparison with the experimental results of Yang, Jones and Karagiozova [IJIE, 2004]



Impact velocity :



4. Complex numerical examples Dynamic buckling of square aluminum tubes

Final results comparison:



Time evolution for the 14.84 m/s impact velocity:





- Advantages of the consistent scheme:
 - Conservation laws and physical consistency are verified for each time step size in the non-linear range
 - Conservation of angular momentum even if numerical dissipation is introduced
- Advantages of the implicit/explicit combined scheme:
 - Reduction of the CPU cost
 - Automatic algorithms
 - No lack of accuracy
 - Remains available after code optimizations



- Drawbacks of the consistent scheme:
 - Mathematical developments needed for each element, material...
 - More complex to implement
- Drawback of the implicit/explicit combined scheme:
 - Implicit and explicit elements must have the same formulation



- Development of a second order accurate EDMC scheme
- Extension to a hyper-elastic model based on an incremental potential (in progress)
- Development of a thermo-mechanical consistent scheme
- Modelization of wind-milling in a turbo-engine